

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM030822\  
 Data File : BM034160.D  
 Acq On : 08 Mar 2022 22:18  
 Operator : CG/JU  
 Sample : N1733-10  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBRS5

Quant Time: Mar 09 00:42:48 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM030822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 08 13:09:16 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.989  | 152  | 152488   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.807 | 136  | 581599   | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10      | 14.630 | 164  | 369164   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.365 | 188  | 724458   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12          | 21.541 | 240  | 681247   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12          | 23.983 | 264  | 762525   | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.372  | 96  | 26590   | 8.020  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.790  | 84  | 313034  | 32.435 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.142  | 99  | 457732  | 38.309 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.313  | 67  | 257780  | 40.084 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.519  | 132 | 404184  | 40.809 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.695  | 113 | 381301  | 38.086 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.154  | 128 | 196088  | 43.514 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.878  | 143 | 212060  | 43.327 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.419 | 165 | 411112  | 40.731 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.936 | 131 | 521794  | 37.838 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 14.036 | 166 | 1164602 | 41.573 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.324 | 160 | 1495140 | 42.681 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.801 | 143 | 160554  | 30.802 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.618 | 176 | 1025610 | 41.137 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.724 | 200 | 142473  | 30.474 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.465 | 188 | 1514504 | 42.439 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.753 | 212 | 1651919 | 44.114 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.830 | 264 | 1725990 | 42.670 | ng/u1 | 0.00 |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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